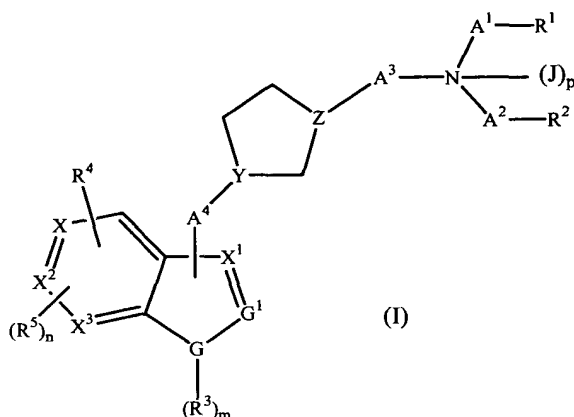


What is claimed is:

1. A compound of Formula (I)

5



or pharmaceutically acceptable salt or solvate thereof

wherein

A^1 and A^2 are each independently C_{1-4} alkylene or a bond;

10 A^3 is a bond, C_{1-4} alkylene or C_{1-4} alkylidene;

A^4 is C_{1-4} alkylene or a bond and is attached to X, X^1 or X^2 ;

X, X^1 , X^2 and X^3 are independently C or CH;

J is C_{1-4} alkyl;

p is 0 or 1;

15 R^1 and R^2 are independently H, C_{1-3} alkyl, C_{3-6} cycloalkyl, phenyl, -O-phenyl, -N(H)C(O)O- C_{1-4} alkyl or C_{1-4} alkyl-N(H)C(O)O-;

said C_{3-6} cycloalkyl, phenyl or O-phenyl being
independently and optionally substituted with
 C_{1-4} alkyl, C_{1-3} alkoxy, indolyl or halo;

20 wherein said indolyl is optionally
substituted by halo or cyano;

or are independently selected from the group of heterocyclic moieties consisting of thienyl, furanyl, pyrrolyl, pyrrolinyl, pyrrolidinyl, imidazolyl, imidazoliny, imidazolidinyl, pyrazolyl, pyrazoliny, pyrazolidinyl, pyridyl, pyrimidinyl, piperidinyl, piperazinyl, morpholino, adamantyl, indolyl, isoindolyl, indoliny, quinoliny, dihydroquinoliny, tetrahydroquinoliny, isoquinoliny, dihydroisoquinoliny and tetrahydroisoquinoliny, wherein said heterocyclic moieties are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy or cyano;

or wherein -A¹-R¹ and -A²-R² together with the nitrogen to which they are attached form pyrrolyl, pyrroliny, pyrrolidinyl, imidazolyl, imidazoliny, imidazolidinyl, pyrazolyl, pyrazoliny, pyrazolidinyl, pyridyl, pyrimidinyl, piperidinyl, piperazinyl, morpholino, indolyl, isoindolyl, indoliny, isoindoliny, quinoliny, dihydroquinoliny, tetrahydroquinoliny, isoquinoliny, dihydroisoquinoliny or tetrahydroisoquinoliny and are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy, cyano or benzyl;

R³ is H or C₁₋₄alkyl;

m is 0 or 1;

R⁴ and R⁵ are independently hydrogen, cyano, halo, nitro, C₁₋₃alkyl or C₁₋₃perfluoroalkyl;

wherein said R⁴ or R⁵ may be independently attached to G¹, X, X¹, X² or X³;

n is 0 or 1;

G is N, O or S;

G¹ is N, C or CH;

Y is (D)H wherein D is C; and

Z is (E)H wherein E is C;

provided that

5 both R⁴ and R⁵ are not attached to the same of said G¹, X, X¹,
X² or X³;

if G is O or S, then m is 0;

if G is N, then m is 1;

10 if R₁ is C₃₋₆cycloalkyl, phenyl or O-phenyl being
independently and optionally substituted with C₁₋₄alkyl,
C₁₋₃alkoxy, indolyl or halo; wherein said indolyl is
optionally substituted by halo or cyano, then R₂ is H or
C₁₋₃alkyl;

15 if R₂ is C₃₋₆cycloalkyl, phenyl or O-phenyl being
independently and optionally substituted with C₁₋₄alkyl,
C₁₋₃alkoxy, indolyl or halo; wherein said indolyl is
optionally substituted by halo or cyano, then R₁ is H or
C₁₋₃alkyl;

20 if -A¹-R¹ and -A²-R² together with the nitrogen to which they
are attached form pyrrolyl, pyrrolinyl, pyrrolidinyl,
imidazolyl, imidazoliny, imidazolidinyl, pyrazolyl,
pyrazolinyl, pyrazolidinyl, pyridyl, pyrimidinyl,
piperidinyl, piperazinyl, morpholino, indolyl,
isoindolyl, indoliny, isoindoliny, quinoliny,
25 dihydroquinoliny, tetrahydroquinoliny, isoquinoliny,
dihydroisoquinoliny or tetrahydroisoquinoliny and are
optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy,
cyano or benzyl, then p is 0;

if R¹ is -N(H)C(O)OC₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or said
heterocyclic moiety wherein said heterocyclic moiety

contains a nitrogen atom and said nitrogen atom is attached to A¹, then A¹ is C₂₋₄alkylene;

if R² is -N(H)C(O)OC₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or said heterocyclic moiety wherein said heterocyclic moiety contains a nitrogen atom and said nitrogen atom is attached to A², then A² is C₂₋₄alkylene;

if R¹ is N(H)C(O)O-C₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or a heterocyclic moiety selected from the group consisting of thienyl, furanyl, pyrrolyl, pyrrolinyl, pyrrolidinyl, imidazolyl, imidazoliny, imidazolidinyl, pyrazolyl, pyrazolinyl, pyrazolidinyl, pyridyl, pyrimidinyl, piperidinyl, piperazinyl, morpholino, adamantyl, indolyl, isoindolyl, indolinyl, quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl, dihydroisoquinolinyl and tetrahydroisoquinolinyl, wherein said heterocyclic moieties are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy or cyano, then R² is H or C₁₋₃alkyl;

if R² is -N(H)C(O)O-C₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or a heterocyclic moiety selected from the group consisting of thienyl, furanyl, pyrrolyl, pyrrolinyl, pyrrolidinyl, imidazolyl, imidazoliny, imidazolidinyl, pyrazolyl, pyrazolinyl, pyrazolidinyl, pyridyl, pyrimidinyl, piperidinyl, piperazinyl, morpholino, adamantyl, indolyl, isoindolyl, indolinyl, quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl, dihydroisoquinolinyl and tetrahydroisoquinolinyl, wherein said heterocyclic moieties are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy or cyano, then R¹ is H or C₁₋₃alkyl;

if R⁴ or R⁵ are attached to G¹, then G¹ is C;

if A⁴, R⁴ or R⁵ are attached to X, then X is C;
 if A⁴, R⁴ or R⁵ are attached to X¹, then X¹ is C;
 if A⁴, R⁴ or R⁵ are attached to X², then X² is C;
 if R⁴ or R⁵ are attached to X³, then X³ is C.

5

2. A compound according to claim 1 wherein p is 0.
3. A compound according to claim 1 wherein G is N and G¹ is CH.
4. A compound according to claim 1 wherein G is S and G¹ is CH.
5. A compound according to claim 1 wherein G is N and G¹ is N.
- 10 6. A compound according to claim 1 wherein G is S and G¹ is N.
7. A compound according to claim 1 wherein G is O and G¹ is N.
8. A compound according to claim 1 wherein R¹ is methyl and R² is methyl.
9. A compound according to claim 1 wherein R¹ is H and R² is C₃₋₆cycloalkyl
 wherein said C₃₋₆cycloalkyl is substituted with indolyl and wherein said indolyl is
 15 optionally substituted by halo or cyano.
10. A compound according to claim 1 wherein A¹ is a bond, R¹ is methyl, A² is a
 bond and R² is methyl.
11. A compound according to claim 1 wherein R¹ and R² are independently H, C₁₋₃alkyl,
 C₃₋₆cycloalkyl, phenyl, -O-phenyl, -N(H)C(O)O-C₁₋₄alkyl or C₁₋₄alkyl-
 20 N(H)C(O)O-; said C₃₋₆cycloalkyl, phenyl or O-phenyl being independently and
 optionally substituted with C₁₋₄alkyl, C₁₋₃alkoxy or halo.
12. A compound according to claim 1 wherein R¹ and R² are independently H, C₁₋₃alkyl,
 phenyl, said phenyl being independently and optionally substituted with
 C₁₋₄alkyl, C₁₋₃alkoxy or halo.
- 25 13. A compound according to claim 1 wherein R¹ and R² are independently H or
 unsubstituted C₁₋₃alkyl or phenyl.
14. A compound according to claim 1 wherein R¹ and R² are independently H or
 unsubstituted C₁₋₃alkyl or phenyl and A¹ and A² are independently C₁₋₄alkylene.
15. A compound according to claim 1 wherein -A¹-R¹ and -A²-R² together with the
 30 nitrogen to which they are attached form pyrrolyl, pyrrolinyl, pyrrolidinyl,
 piperidinyl, piperazinyl, morpholino, indolyl, isoindolyl, indolinyl, isoindolinyl,
 quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl,

- dihydroisoquinolinyl or tetrahydroisoquinolinyl and are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy, cyano or benzyl.
16. A compound according to claim 1 wherein -A¹-R¹ and -A²-R² together with the nitrogen to which they are attached form unsubstituted pyrrolyl, pyrrolinyl, 5 pyrrolidinyl, piperidinyl, piperazinyl, morpholino, indolyl, isoindolyl, indolinyl, isoindolinyl, quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl, dihydroisoquinolinyl or tetrahydroisoquinolinyl.
17. A compound according to claim 1 wherein -A¹-R¹ and -A²-R² together with the nitrogen to which they are attached form unsubstituted pyrrolidinyl, piperidinyl, 10 morpholino or isoindolinyl.
18. A compound according to claim 1 wherein R³ is H and m is 1.
19. A compound according to claim 1 wherein n is 0.
20. A compound according to claim 1 wherein R⁴ and R⁵ are halo.
21. A compound according to claim 1 wherein R⁴ is C₁₋₃alkyl and is attached to G¹.
- 15 22. A compound according to claim 1 wherein R⁴ is C₁₋₃perfluoroalkyl and is attached to G¹.
23. A compound according to claim 1 wherein R⁴ is hydrogen.
24. A compound according to claim 1 wherein R⁴ is fluoro.
25. A compound according to claim 1 wherein R⁴ is cyano.
- 20 26. A compound according to claim 1 wherein R⁴ is cyano or fluoro.
27. A compound according to claim 1 wherein R⁴ and R⁵ are each fluoro.
28. A compound according to claim 1 wherein the hydrogen atom attached to D is in the *trans* configuration to the hydrogen atom attached to E.
29. A compound according to claim 1 wherein the hydrogen atom attached to D is in the *cis* configuration to the hydrogen atom attached to E.
- 25 30. A compound according to claim 1 wherein D in relation to the four moieties to which it is attached has an absolute configuration of S; E in relation to the four moieties to which it is attached has an absolute configuration of S.
31. A compound according to claim 1 wherein D in relation to the four moieties to which it is attached has an absolute configuration of S; E in relation to the four 30 moieties to which it is attached has an absolute configuration of R.

32. A compound according to claim 1 wherein D in relation to the four moieties to which it is attached has an absolute configuration of R; E in relation to the four moieties to which it is attached has an absolute configuration of S.
33. A compound according to claim 1 wherein D in relation to the four moieties to which it is attached has an absolute configuration of R; E in relation to the four moieties to which it is attached has an absolute configuration of R.
34. A compound according to claim 1 wherein A³ is a bond.
35. A compound according to claim 1 wherein A³ is C₁₋₄alkylene.
36. A compound according to claim 1 wherein A³ is C₁₋₄alkylidene.
37. A compound according to claim 1 wherein A³ is methylene.
38. A compound according to claim 1 wherein A⁴ is a bond.
39. A compound according to claim 1 wherein A⁴ is methylene.
40. A compound according to claim 1 wherein A⁴ is attached X¹.
41. A compound according to claim 1 wherein A⁴ is attached X.
42. A compound according to claim 1 wherein R⁴ is attached X.
43. A compound according to claim 1 wherein
- A¹ and A² are each independently C₁₋₄alkylene or a bond;
- A³ is a bond;
- A⁴ is a bond and is attached to X¹;
- X and X¹ are each C;
- X² and X³ are each CH;
- p is 0;
- R¹ and R² are independently H, C₁₋₃alkyl, C₃₋₆cycloalkyl, phenyl, -O-phenyl, -N(H)C(O)O-C₁₋₄alkyl or C₁₋₄alkyl-N(H)C(O)O-;
- said C₃₋₆cycloalkyl, phenyl or O-phenyl being independently and optionally substituted with C₁₋₄alkyl, C₁₋₃alkoxy or halo;
- or are independently selected from the group of heterocyclic moieties consisting of thienyl, furanyl, pyrrolyl, pyrrolinyl, pyrrolidinyl, imidazolyl, imidazoliny,

5 imidazolidinyl, pyrazolyl, pyrazolinyl, pyrazolidinyl,
pyridyl, pyrimidinyl, piperidinyl, piperazinyl,
morpholino, adamantyl, indolyl, isoindolyl, indolinyl,
quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl,
isoquinolinyl, dihydroisoquinolinyl and
tetrahydroisoquinolinyl, wherein said heterocyclic
moieties are optionally substituted with halo, C₁₋₄alkyl,
C₁₋₄alkoxy or cyano;

10 or wherein -A¹-R¹ and -A²-R² together with the nitrogen to
which they are attached form pyrrolyl, pyrrolinyl,
pyrrolidinyl, imidazolyl, imidazolinyl, imidazolidinyl,
pyrazolyl, pyrazolinyl, pyrazolidinyl, pyridyl,
pyrimidinyl, piperidinyl, piperazinyl, morpholino,
15 indolyl, isoindolyl, indolinyl, isoindolinyl, quinolinyl,
dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl,
dihydroisoquinolinyl or tetrahydroisoquinolinyl and are
optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy,
cyano or benzyl;

R³ is H;

20 m is 1;

R⁴ is hydrogen, cyano, halo, nitro, C₁₋₃alkyl or C₁₋₃perfluoroalkyl and
is attached to X;

n is 0;

G is N;

25 G¹ is CH;

Y is (D)H wherein D is C; and

Z is (E)H wherein E is C;

provided that

- if R¹ is -N(H)C(O)OC₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or said heterocyclic moiety wherein said heterocyclic moiety contains a nitrogen atom and said nitrogen atom is attached to A¹, then A¹ is C₂₋₄alkylene;
- 5 if R² is -N(H)C(O)OC₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or said heterocyclic moiety wherein said heterocyclic moiety contains a nitrogen atom and said nitrogen atom is attached to A², then A² is C₂₋₄alkylene;
- 10 if R¹ is N(H)C(O)O-C₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or a heterocyclic moiety selected from the group consisting of thienyl, furanyl, pyrrolyl, pyrrolinyl, pyrrolidinyl, imidazolyl, imidazolinyl, imidazolidinyl, pyrazolyl, pyrazolinyl, pyrazolidinyl, pyridyl, pyrimidinyl, piperidinyl, piperazinyl, morpholino, adamantyl,
- 15 indolyl, isoindolyl, indolinyl, quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl, dihydroisoquinolinyl and tetrahydroisoquinolinyl, wherein said heterocyclic moieties are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy or cyano,
- 20 then R² is H or C₁₋₃alkyl; and
- if R² is -N(H)C(O)O-C₁₋₄alkyl, C₁₋₄alkyl-N(H)C(O)O- or a heterocyclic moiety selected from the group consisting of thienyl, furanyl, pyrrolyl, pyrrolinyl, pyrrolidinyl, imidazolyl, imidazolinyl, imidazolidinyl, pyrazolyl, pyrazolinyl, pyrazolidinyl, pyridyl, pyrimidinyl,
- 25 piperidinyl, piperazinyl, morpholino, adamantyl, indolyl, isoindolyl, indolinyl, quinolinyl, dihydroquinolinyl, tetrahydroquinolinyl, isoquinolinyl, dihydroisoquinolinyl and tetrahydroisoquinolinyl, wherein said heterocyclic moieties are optionally substituted with halo, C₁₋₄alkyl, C₁₋₄alkoxy or cyano,
- 30 then R¹ is H or C₁₋₃alkyl.

44. A pharmaceutically acceptable formulation comprising a compound according to claim 1.
45. A method of treating depression, attention deficit hyperactivity disorder,
5 obsessive-compulsive disorder, post-traumatic stress disorder, substance abuse disorders and sexual dysfunction comprising the administration to a human in need thereof an effective amount of a pharmaceutically acceptable formulation comprising a compound according to claim 1.
46. A method of treating sexual dysfunction comprising the administration to a
10 human in need thereof an effective amount of a pharmaceutically acceptable formulation comprising a compound according to claim 1.
47. A method of treating premature ejaculation comprising the administration to a human in need thereof an effective amount of a pharmaceutically acceptable formulation comprising a compound according to claim 1.
- 15 48. A compound or pharmaceutically acceptable salt or solvate thereof selected from the group consisting of
- 3-(3-methylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-(3-ethylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-(3-dimethylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 20 3-[3-(ethyl-methyl-amino)-cyclopentyl]-1*H*-indole-5-carbonitrile;
 - 3-(3-diethylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-(3-pyrrolidin-1-yl-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-[3-(1,3-dihydro-isoindol-2-yl)-cyclopentyl]-1*H*-indole-5-carbonitrile;
 - 3-[3-(3,4-dihydro-1*H*-isoquinolin-2-yl)-cyclopentyl]-1*H*-indole-5-carbonitrile;
 - 25 3-(3-penethylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-[3-(methyl-phenethyl-amino)-cyclopentyl]-1*H*-indole-5-carbonitrile;
 - 3-(3-morpholin-4-yl-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-[3-(benzyl-methyl-amino)-cyclopentyl]-1*H*-indole-5-carbonitrile;
 - 3-(3-benzylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 30 3-(3-piperidin-1-yl-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-(3-dipropylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 - 3-(3-propylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;

- 1-methyl-3-(3-methylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 3-(3-ethylamino-cyclopentyl)-1-methyl-1*H*-indole-5-carbonitrile;
 3-(3-benzylamino-cyclopentyl)-1-methyl-1*H*-indole-5-carbonitrile;
 1-methyl-3-(3-phenethylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 5 3-(3-dimethylamino-cyclopentyl)-1-methyl-1*H*-indole-5-carbonitrile;
 3-[3-(ethyl-methyl-amino)-cyclopentyl]-1-methyl-1*H*-indole-5-carbonitrile;
 3-(3-diethylamino-cyclopentyl)-1-methyl-1*H*-indole-5-carbonitrile;
 1-methyl-3-(3-pyrrolidin-1-yl-cyclopentyl)-1*H*-indole-5-carbonitrile;
 1-methyl-3-(3-piperidin-1-yl-cyclopentyl)-1*H*-indole-5-carbonitrile;
 10 1-methyl-3-(3-morpholin-4-yl-cyclopentyl)-1*H*-indole-5-carbonitrile;
 3-[3-(benzyl-methyl-amino)-cyclopentyl]-1-methyl-1*H*-indole-5-carbonitrile;
 1-methyl-3-[3-(methyl-phenethyl-amino)-cyclopentyl]-1*H*-indole-5-carbonitrile;
 1-methyl-3-(3-propylamino-cyclopentyl)-1*H*-indole-5-carbonitrile;
 3-(3-dipropylamino-cyclopentyl)-1-methyl-1*H*-indole-5-carbonitrile;
 15 3-[3-(benzyl-methyl-amino)-cyclopentyl]-1-ethyl-1*H*-indole-5-carbonitrile;
 3-(3-dimethylamino-cyclopentyl)-1-ethyl-1*H*-indole-5-carbonitrile;
 3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl]-dimethyl-amine;
 ethyl-[3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl]-methyl-amine;
 diethyl-[3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl]-amine;
 20 5-fluoro-3-(3-pyrrolidin-1-yl-cyclopentyl)-1*H*-indole;
 3-(4-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(4-bromo-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(5-dihalo-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(5-bromo-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 25 3-(5-iodo-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(6-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(6-chloro-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(6-bromo-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(7-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 30 3-(7-chloro-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 3-(7-bromo-1*H*-indol-3-yl)-cyclopentyl-dimethyl-amine;
 (1*S*,3*R*)-3-(3-dimethylaminocyclopentyl)-1*H*-indole-5-carbonitrile;
 (1*S*,3*S*)-3-(3-dimethylaminocyclopentyl)-1*H*-indole-5-carbonitrile;

- (1*R*,3*S*)-3-(3-dimethylaminocyclopentyl)-1*H*-indole-5-carbonitrile;
(1*R*,3*R*)-3-(3-dimethylaminocyclopentyl)-1*H*-indole-5-carbonitrile;
(1*S*,3*S*)-3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethylamine;
(1*R*,3*S*)-3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethylamine;
5 (1*R*,3*R*)-3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethylamine;
(1*S*,3*R*)-3-(5-fluoro-1*H*-indol-3-yl)-cyclopentyl-dimethylamine;
(1*S*,3*R*)-3-(3-dimethylamino-cyclopentyl)-1-ethyl-1*H*-indole-5-carbonitrile;
(1*S*,3*S*)-3-(3-dimethylamino-cyclopentyl)-1-ethyl-1*H*-indole-5-carbonitrile;
(1*R*,3*S*)-3-(3-dimethylamino-cyclopentyl)-1-ethyl-1*H*-indole-5-carbonitrile;
10 (1*R*,3*R*)-3-(3-dimethylamino-cyclopentyl)-1-ethyl-1*H*-indole-5-carbonitrile;
(1*S*)-3-(3-amino-cyclopentyl)-1*H*-indole-5-carbonitrile; and
(3*S*,3'*S*)-bis-(3-(5-cyano-1*H*-indol-3-yl)cyclopentyl)amine.